



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:53 PM UTC

PDB ID : 3QWY / pdb_00003qwy
Title : CED-2
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Deposited on : 2011-02-28
Resolution : 2.52 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

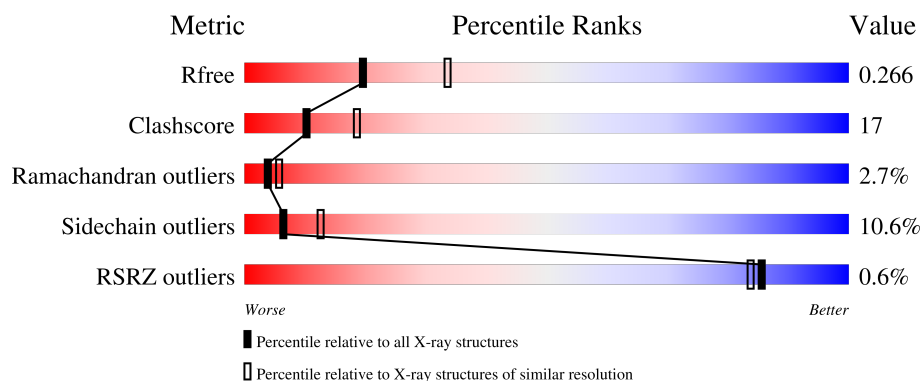
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

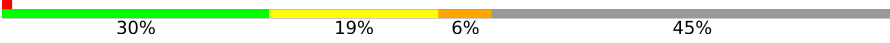
The reported resolution of this entry is 2.52 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7383 (2.54-2.50)
Clashscore	190562	8079 (2.54-2.50)
Ramachandran outliers	187476	7944 (2.54-2.50)
Sidechain outliers	187428	7946 (2.54-2.50)
RSRZ outliers	180081	7387 (2.54-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	308	 30% 20% . . 45%
1	B	308	 % 30% 19% 6% 45%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	B	280	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2830 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Cell death abnormality protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1345	858	228	255	4			
1	B	170	Total	C	N	O	S	0	0	0
			1346	857	229	256	4			

There are 58 discrepancies between the modelled and reference sequences:

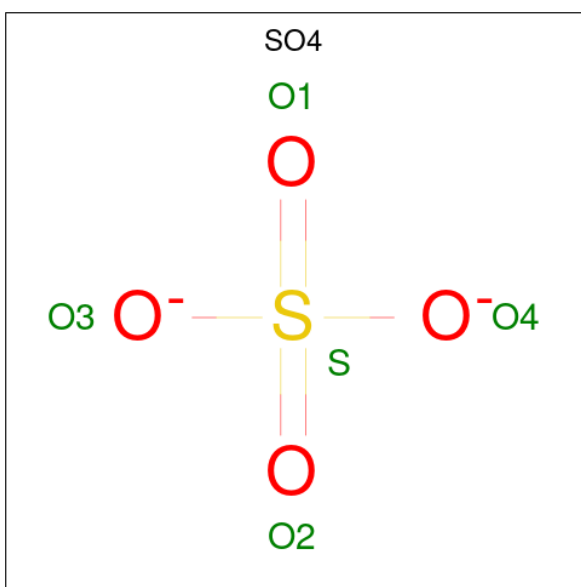
Chain	Residue	Modelled	Actual	Comment	Reference
A	-28	MET	-	expression tag	UNP Q9NHC3
A	-27	GLY	-	expression tag	UNP Q9NHC3
A	-26	SER	-	expression tag	UNP Q9NHC3
A	-25	SER	-	expression tag	UNP Q9NHC3
A	-24	HIS	-	expression tag	UNP Q9NHC3
A	-23	HIS	-	expression tag	UNP Q9NHC3
A	-22	HIS	-	expression tag	UNP Q9NHC3
A	-21	HIS	-	expression tag	UNP Q9NHC3
A	-20	HIS	-	expression tag	UNP Q9NHC3
A	-19	HIS	-	expression tag	UNP Q9NHC3
A	-18	SER	-	expression tag	UNP Q9NHC3
A	-17	SER	-	expression tag	UNP Q9NHC3
A	-16	GLY	-	expression tag	UNP Q9NHC3
A	-15	LEU	-	expression tag	UNP Q9NHC3
A	-14	VAL	-	expression tag	UNP Q9NHC3
A	-13	PRO	-	expression tag	UNP Q9NHC3
A	-12	ARG	-	expression tag	UNP Q9NHC3
A	-11	GLY	-	expression tag	UNP Q9NHC3
A	-10	SER	-	expression tag	UNP Q9NHC3
A	-9	HIS	-	expression tag	UNP Q9NHC3
A	-8	MET	-	expression tag	UNP Q9NHC3
A	-7	ALA	-	expression tag	UNP Q9NHC3
A	-6	SER	-	expression tag	UNP Q9NHC3
A	-5	MET	-	expression tag	UNP Q9NHC3
A	-4	THR	-	expression tag	UNP Q9NHC3

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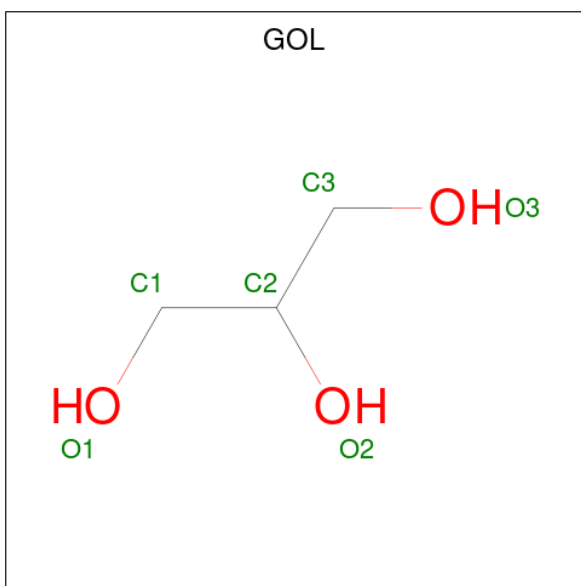
Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q9NHC3
A	-2	GLY	-	expression tag	UNP Q9NHC3
A	-1	GLN	-	expression tag	UNP Q9NHC3
A	0	GLN	-	expression tag	UNP Q9NHC3
B	-28	MET	-	expression tag	UNP Q9NHC3
B	-27	GLY	-	expression tag	UNP Q9NHC3
B	-26	SER	-	expression tag	UNP Q9NHC3
B	-25	SER	-	expression tag	UNP Q9NHC3
B	-24	HIS	-	expression tag	UNP Q9NHC3
B	-23	HIS	-	expression tag	UNP Q9NHC3
B	-22	HIS	-	expression tag	UNP Q9NHC3
B	-21	HIS	-	expression tag	UNP Q9NHC3
B	-20	HIS	-	expression tag	UNP Q9NHC3
B	-19	HIS	-	expression tag	UNP Q9NHC3
B	-18	SER	-	expression tag	UNP Q9NHC3
B	-17	SER	-	expression tag	UNP Q9NHC3
B	-16	GLY	-	expression tag	UNP Q9NHC3
B	-15	LEU	-	expression tag	UNP Q9NHC3
B	-14	VAL	-	expression tag	UNP Q9NHC3
B	-13	PRO	-	expression tag	UNP Q9NHC3
B	-12	ARG	-	expression tag	UNP Q9NHC3
B	-11	GLY	-	expression tag	UNP Q9NHC3
B	-10	SER	-	expression tag	UNP Q9NHC3
B	-9	HIS	-	expression tag	UNP Q9NHC3
B	-8	MET	-	expression tag	UNP Q9NHC3
B	-7	ALA	-	expression tag	UNP Q9NHC3
B	-6	SER	-	expression tag	UNP Q9NHC3
B	-5	MET	-	expression tag	UNP Q9NHC3
B	-4	THR	-	expression tag	UNP Q9NHC3
B	-3	GLY	-	expression tag	UNP Q9NHC3
B	-2	GLY	-	expression tag	UNP Q9NHC3
B	-1	GLN	-	expression tag	UNP Q9NHC3
B	0	GLN	-	expression tag	UNP Q9NHC3

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	91	Total 91	O 91	0	0
4	B	37	Total 37	O 37	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	105.21Å 97.15Å 71.10Å 90.00° 127.14° 90.00°	Depositor
Resolution (Å)	48.58 – 2.52 48.58 – 2.52	Depositor EDS
% Data completeness (in resolution range)	98.1 (48.58-2.52) 98.1 (48.58-2.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.204 , 0.268 0.201 , 0.266	Depositor DCC
R_{free} test set	965 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	57.7	Xtriage
Anisotropy	0.090	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2830	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.97	24/1378 (1.7%)	1.74	27/1867 (1.4%)
1	B	1.85	19/1378 (1.4%)	1.66	23/1866 (1.2%)
All	All	1.91	43/2756 (1.6%)	1.70	50/3733 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	135	LEU	CA-C	8.66	1.63	1.52
1	A	56	ALA	CA-C	8.52	1.63	1.52
1	B	44	SER	N-CA	8.15	1.56	1.46
1	A	167	PRO	CA-C	7.68	1.59	1.52
1	B	121	VAL	CA-CB	7.51	1.65	1.55
1	A	128	THR	CA-CB	7.26	1.64	1.53
1	B	117	ILE	CA-CB	-7.19	1.45	1.54
1	B	163	THR	CA-CB	7.15	1.64	1.53
1	A	61	ALA	CA-CB	6.83	1.64	1.53
1	A	72	PRO	CA-C	6.78	1.60	1.52
1	A	135	LEU	C-O	-6.68	1.17	1.24
1	A	112	ALA	CA-CB	-6.57	1.43	1.53
1	B	158	ASN	CA-C	6.49	1.60	1.53
1	A	104	LEU	C-O	-6.41	1.16	1.24
1	A	68	GLU	CA-C	-6.22	1.45	1.52
1	B	85	ALA	CA-CB	6.18	1.64	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	16	PHE	CA-C	6.18	1.60	1.52
1	A	125	PHE	N-CA	6.18	1.53	1.46
1	A	85	ALA	CA-CB	6.06	1.63	1.53
1	B	90	PRO	N-CA	5.97	1.55	1.47
1	A	21	ARG	C-O	-5.91	1.17	1.24
1	A	67	ILE	CA-CB	-5.86	1.47	1.54
1	B	31	PRO	N-CA	5.83	1.55	1.47
1	B	166	VAL	C-N	5.81	1.39	1.33
1	A	55	GLU	N-CA	5.76	1.53	1.45
1	A	32	GLN	CG-CD	5.70	1.66	1.52
1	A	18	GLY	N-CA	-5.56	1.35	1.44
1	A	136	PRO	N-CA	5.52	1.53	1.47
1	A	56	ALA	C-O	5.52	1.30	1.23
1	B	125	PHE	N-CA	5.42	1.52	1.46
1	A	32	GLN	CD-OE1	5.40	1.33	1.23
1	A	53	VAL	C-O	5.38	1.29	1.24
1	A	9	PHE	N-CA	5.34	1.53	1.46
1	B	156	ALA	C-O	5.33	1.30	1.23
1	B	163	THR	N-CA	5.33	1.52	1.45
1	B	85	ALA	CA-C	5.31	1.59	1.52
1	A	130	GLU	CD-OE1	5.30	1.35	1.25
1	B	7	ASP	CA-C	5.18	1.58	1.52
1	A	62	VAL	C-O	5.12	1.29	1.24
1	B	67	ILE	CA-CB	5.08	1.60	1.53
1	B	44	SER	C-O	-5.05	1.17	1.24
1	A	85	ALA	N-CA	5.02	1.52	1.46
1	B	45	ARG	CA-C	5.00	1.59	1.53

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	13	SER	N-CA-C	10.20	124.83	112.38
1	A	117	ILE	CB-CA-C	-9.58	99.03	110.91
1	A	151	GLN	N-CA-C	9.33	122.61	111.33
1	A	14	PHE	N-CA-C	-9.06	101.19	112.88
1	A	25	HIS	N-CA-C	-8.89	101.67	111.36
1	B	44	SER	N-CA-C	-8.34	103.25	113.50
1	B	16	PHE	CA-C-N	-8.28	111.36	120.14
1	B	16	PHE	C-N-CA	-8.28	111.36	120.14
1	A	86	ASN	N-CA-C	-7.92	93.94	110.80
1	B	152	ASP	N-CA-C	7.52	122.46	113.28
1	A	71	GLU	N-CA-C	-7.42	100.08	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	173	ILE	CA-C-N	7.23	134.71	121.70
1	B	173	ILE	C-N-CA	7.23	134.71	121.70
1	A	139	GLN	N-CA-C	-7.07	98.51	109.76
1	A	30	GLU	CA-C-N	-6.71	112.86	119.64
1	A	30	GLU	C-N-CA	-6.71	112.86	119.64
1	A	115	LYS	CA-C-N	-6.69	113.75	120.31
1	A	115	LYS	C-N-CA	-6.69	113.75	120.31
1	B	89	PHE	CA-C-N	-6.69	112.74	119.56
1	B	89	PHE	C-N-CA	-6.69	112.74	119.56
1	B	92	ILE	O-C-N	6.48	124.56	120.42
1	A	45	ARG	N-CA-C	-6.32	100.46	109.24
1	A	10	GLU	CB-CA-C	-6.27	99.63	110.72
1	A	9	PHE	N-CA-C	6.23	120.14	112.54
1	A	113	TYR	N-CA-C	6.00	119.30	110.30
1	A	76	GLY	N-CA-C	-5.92	107.15	115.32
1	B	74	GLU	N-CA-C	5.88	119.28	111.75
1	A	79	ALA	N-CA-C	-5.80	98.44	110.80
1	B	167	PRO	CB-CA-C	-5.79	106.13	111.40
1	B	71	GLU	CB-CA-C	5.78	116.71	108.68
1	B	113	TYR	N-CA-C	-5.76	101.16	110.32
1	A	105	THR	N-CA-C	-5.62	104.43	112.13
1	A	87	GLN	N-CA-C	5.61	117.36	108.67
1	A	130	GLU	O-C-N	-5.60	116.42	122.19
1	A	170	TYR	N-CA-C	-5.59	105.77	112.59
1	B	5	GLY	N-CA-C	-5.54	97.22	113.30
1	B	161	GLY	N-CA-C	5.53	126.29	113.18
1	B	164	GLY	CA-C-O	-5.51	116.20	121.83
1	B	45	ARG	CG-CD-NE	-5.42	100.08	112.00
1	B	129	GLY	N-CA-C	5.40	119.01	112.48
1	A	90	PRO	N-CA-C	-5.31	107.22	113.86
1	A	136	PRO	N-CA-C	5.30	119.64	111.21
1	B	120	VAL	CA-C-N	-5.16	116.09	123.17
1	B	120	VAL	C-N-CA	-5.16	116.09	123.17
1	B	23	GLU	CB-CA-C	-5.15	102.80	110.88
1	A	116	PRO	CA-C-N	5.12	128.98	122.37
1	A	116	PRO	C-N-CA	5.12	128.98	122.37
1	A	10	GLU	N-CA-CB	5.04	117.88	110.57
1	B	51	LEU	N-CA-C	-5.02	100.23	108.41
1	B	103	VAL	N-CA-C	5.01	117.04	109.12

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	78	ALA	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1345	0	1307	45	0
1	B	1346	0	1305	46	0
2	A	5	0	0	0	0
3	B	6	0	8	3	0
4	A	91	0	0	6	0
4	B	37	0	0	3	0
All	All	2830	0	2620	91	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 17.

All (91) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LYS:H	1:A:78:ALA:HB2	1.07	1.12
1:B:21:ARG:HH11	3:B:280:GOL:H32	1.16	1.07
1:A:73:LYS:H	1:A:78:ALA:CB	1.73	1.01
1:A:73:LYS:N	1:A:78:ALA:HB2	1.76	0.99
1:A:75:ASP:HB3	1:A:77:THR:OG1	1.60	0.98
1:B:25:HIS:HD2	1:B:64:HIS:CE1	1.85	0.94
1:B:25:HIS:HD2	1:B:64:HIS:HE1	0.95	0.94
1:B:25:HIS:CD2	1:B:64:HIS:HE1	1.85	0.92
1:A:41:ARG:HH21	1:A:64:HIS:CD2	1.89	0.91
1:A:11:TRP:HA	1:A:14:PHE:CE2	2.11	0.86
1:A:74:GLU:O	1:A:76:GLY:N	2.11	0.83
1:A:41:ARG:HH21	1:A:64:HIS:HD2	1.25	0.80
1:A:22:GLU:HG3	4:A:283:HOH:O	1.82	0.78
1:A:74:GLU:O	1:A:75:ASP:C	2.27	0.77
1:B:174:GLN:O	4:B:283:HOH:O	2.08	0.72
1:A:119:GLU:HB2	1:A:145:ILE:HB	1.71	0.70
1:A:98:HIS:HE1	4:A:338:HOH:O	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:ARG:NH1	3:B:280:GOL:H32	2.01	0.64
1:B:41:ARG:HH21	1:B:64:HIS:HD2	1.45	0.63
1:B:21:ARG:O	1:B:21:ARG:HG2	1.97	0.63
1:A:84:ILE:HG22	1:A:95:LEU:HD11	1.81	0.62
1:B:174:GLN:C	4:B:283:HOH:O	2.43	0.62
1:B:20:SER:OG	1:B:23:GLU:HG3	2.00	0.61
1:A:73:LYS:N	1:A:78:ALA:CB	2.51	0.59
1:A:75:ASP:CB	1:A:77:THR:OG1	2.43	0.59
1:B:54:ARG:NH2	1:B:57:ASP:OD1	2.35	0.58
1:B:146:LEU:N	1:B:146:LEU:HD23	2.19	0.58
1:A:78:ALA:O	1:A:79:ALA:CB	2.52	0.58
1:B:45:ARG:CZ	1:B:45:ARG:HA	2.34	0.58
1:B:32:GLN:HG2	1:B:33:VAL:H	1.69	0.57
1:A:10:GLU:HG3	4:A:352:HOH:O	2.03	0.57
1:B:40:MET:HE1	1:B:92:ILE:HD13	1.87	0.56
1:A:41:ARG:NH2	1:A:64:HIS:CD2	2.69	0.55
1:B:30:GLU:HB3	1:B:33:VAL:HG23	1.88	0.55
1:A:74:GLU:C	1:A:76:GLY:N	2.65	0.55
1:B:173:ILE:HG22	1:B:174:GLN:HG2	1.89	0.54
1:B:21:ARG:HH11	3:B:280:GOL:C3	2.04	0.54
1:B:173:ILE:O	1:B:174:GLN:HG3	2.08	0.54
1:B:117:ILE:HG12	1:B:144:GLU:HG2	1.90	0.54
1:B:119:GLU:HA	1:B:119:GLU:OE1	2.08	0.53
1:B:42:ASP:HB2	4:B:309:HOH:O	2.08	0.52
1:A:121:VAL:HB	1:A:171:VAL:HG13	1.90	0.52
1:A:15:TYR:OH	1:A:42:ASP:OD1	2.27	0.51
1:B:91:ASP:HB2	1:B:93:PRO:HD2	1.92	0.51
1:B:41:ARG:HH21	1:B:64:HIS:CD2	2.29	0.50
1:A:11:TRP:HA	1:A:14:PHE:CD2	2.46	0.50
1:A:156:ALA:O	1:A:163:THR:HA	2.12	0.49
1:B:30:GLU:O	1:B:54:ARG:NH1	2.45	0.49
1:A:148:LYS:HG2	1:A:154:TRP:CD1	2.48	0.49
1:A:55:GLU:HG2	1:A:107:ALA:HB3	1.94	0.48
1:B:32:GLN:HG2	1:B:33:VAL:N	2.29	0.48
1:B:32:GLN:OE1	1:B:32:GLN:N	2.25	0.48
1:B:37:THR:HA	1:B:111:ALA:O	2.14	0.48
1:A:84:ILE:CG2	1:A:95:LEU:HD11	2.44	0.47
1:B:41:ARG:NH2	1:B:64:HIS:HD2	2.09	0.47
1:B:142:ARG:O	1:B:143:LEU:HD23	2.15	0.47
1:A:121:VAL:HG11	1:A:171:VAL:CG1	2.45	0.47
1:B:30:GLU:HG3	1:B:32:GLN:NE2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:MET:HG3	1:A:23:GLU:HB3	1.97	0.46
1:A:52:THR:OG1	1:A:64:HIS:CD2	2.68	0.46
1:A:119:GLU:HA	1:A:119:GLU:OE1	2.15	0.46
1:A:52:THR:OG1	1:A:64:HIS:HD2	1.99	0.46
1:B:65:TYR:CD2	1:B:65:TYR:N	2.83	0.46
1:A:78:ALA:O	1:A:79:ALA:HB3	2.15	0.46
1:B:152:ASP:N	1:B:152:ASP:OD1	2.49	0.45
1:B:55:GLU:HB3	1:B:106:GLU:HB2	1.99	0.45
1:A:84:ILE:HD13	1:A:84:ILE:HG21	1.61	0.45
1:A:148:LYS:HG2	1:A:154:TRP:NE1	2.30	0.45
1:A:84:ILE:O	1:A:85:ALA:HB3	2.17	0.45
1:B:11:TRP:HA	1:B:14:PHE:CE2	2.52	0.45
1:B:158:ASN:C	1:B:158:ASN:OD1	2.60	0.45
1:A:25:HIS:CD2	4:A:321:HOH:O	2.70	0.44
1:A:174:GLN:C	4:A:293:HOH:O	2.59	0.44
1:B:138:GLU:HA	1:B:138:GLU:OE1	2.16	0.44
1:B:31:PRO:HD2	1:B:32:GLN:OE1	2.18	0.43
1:A:121:VAL:CG1	1:A:171:VAL:HG13	2.48	0.43
1:A:25:HIS:HD2	4:A:321:HOH:O	2.02	0.42
1:A:19:MET:HE3	1:A:19:MET:HB2	1.80	0.42
1:B:32:GLN:H	1:B:32:GLN:CD	2.20	0.42
1:A:41:ARG:HH11	1:A:41:ARG:HD2	1.68	0.42
1:B:26:LYS:HG2	1:B:27:LEU:N	2.34	0.42
1:B:115:LYS:HE2	1:B:115:LYS:HB3	1.65	0.42
1:A:30:GLU:OE2	1:A:30:GLU:HA	2.18	0.42
1:A:146:LEU:O	1:A:147:SER:HB2	2.20	0.42
1:A:65:TYR:HB3	1:A:84:ILE:HD11	2.02	0.41
1:B:11:TRP:O	1:B:12:ARG:C	2.64	0.41
1:B:41:ARG:NH2	1:B:64:HIS:CD2	2.88	0.41
1:B:43:SER:HB3	1:B:48:GLU:HB2	2.03	0.41
1:B:26:LYS:O	1:B:27:LEU:C	2.64	0.41
1:B:119:GLU:OE1	1:B:119:GLU:CA	2.67	0.41
1:A:53:VAL:HG12	1:A:55:GLU:HG3	2.01	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	168/308 (54%)	157 (94%)	6 (4%)	5 (3%)	3	4
1	B	168/308 (54%)	152 (90%)	12 (7%)	4 (2%)	4	7
All	All	336/616 (54%)	309 (92%)	18 (5%)	9 (3%)	4	6

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	ASP
1	A	78	ALA
1	A	85	ALA
1	A	86	ASN
1	A	79	ALA
1	B	33	VAL
1	B	160	LEU
1	B	27	LEU
1	B	104	LEU

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	141/257 (55%)	128 (91%)	13 (9%)	8	17
1	B	141/257 (55%)	124 (88%)	17 (12%)	5	9
All	All	282/514 (55%)	252 (89%)	30 (11%)	6	12

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	30	GLU
1	A	33	VAL
1	A	45	ARG
1	A	73	LYS
1	A	77	THR
1	A	100	LYS
1	A	102	ARG
1	A	108	SER
1	A	114	LYS
1	A	130	GLU
1	A	132	GLU
1	A	171	VAL
1	B	26	LYS
1	B	30	GLU
1	B	51	LEU
1	B	58	GLU
1	B	71	GLU
1	B	86	ASN
1	B	87	GLN
1	B	108	SER
1	B	115	LYS
1	B	128	THR
1	B	133	THR
1	B	144	GLU
1	B	146	LEU
1	B	147	SER
1	B	162	THR
1	B	171	VAL
1	B	174	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	64	HIS
1	B	25	HIS
1	B	64	HIS
1	B	86	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	B	280	-	5,5,5	1.08	1 (20%)	5,5,5	3.79	3 (60%)
2	SO4	A	280	-	4,4,4	0.65	0	6,6,6	1.21	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	280	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	280	GOL	C3-C2	-2.09	1.43	1.51

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	280	GOL	O3-C3-C2	-7.00	78.86	110.38
3	B	280	GOL	O2-C2-C3	-3.85	93.24	109.18
3	B	280	GOL	C3-C2-C1	2.07	119.40	111.80

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	280	GOL	O1-C1-C2-C3
3	B	280	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	280	GOL	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	170/308 (55%)	-0.38	0 100 100	36, 47, 66, 79	0
1	B	170/308 (55%)	-0.01	2 (1%) 76 74	39, 57, 80, 98	0
All	All	340/616 (55%)	-0.19	2 (0%) 85 84	36, 51, 76, 98	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	31	PRO	2.3
1	B	85	ALA	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	280	6/6	0.93	0.15	40,44,57,61	0
2	SO4	A	280	5/5	0.99	0.04	47,47,51,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.